



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 11:59 pm BST

PDB ID : 4BEL
Title : BACE2 XAPERONE COMPLEX
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Deposited on : 2013-03-11
Resolution : 1.85 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

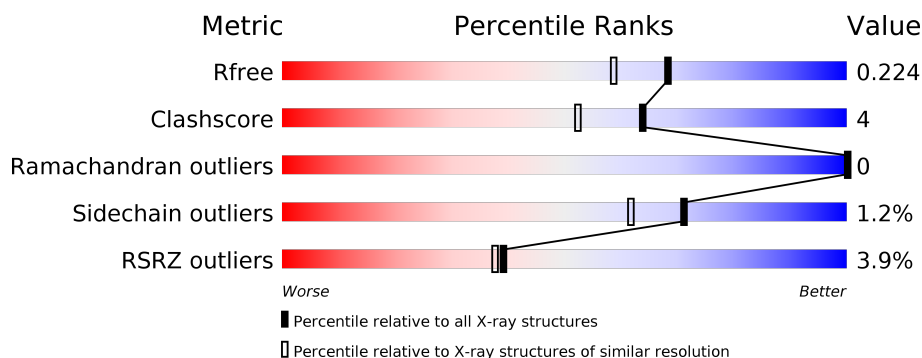
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	386	<div> <div>5%</div> <div> <div></div> <div>85%</div> <div>8%</div> <div>7%</div> </div> </div>
1	B	386	<div> <div>4%</div> <div> <div></div> <div>84%</div> <div>8%</div> <div>8%</div> </div> </div>
2	D	122	<div> <div></div> <div> <div></div> <div>89%</div> <div>..</div> <div>8%</div> </div> </div>
2	E	122	<div> <div></div> <div> <div></div> <div>86%</div> <div>7%</div> <div>7%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8059 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called BETA-SECRETASE 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	358	Total	C	N	O	S	0	6	0
			2821	1816	452	541	12			
1	B	357	Total	C	N	O	S	0	9	0
			2834	1821	454	547	12			

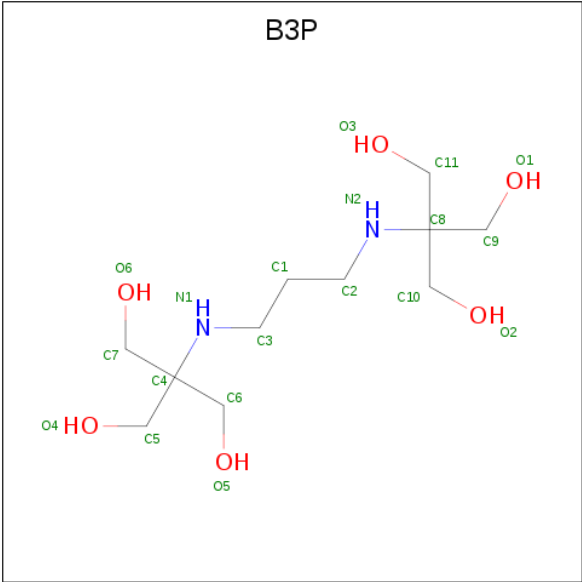
- Molecule 2 is a protein called XA4813.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	112	Total	C	N	O	S	0	4	0
			860	527	155	171	7			
2	E	113	Total	C	N	O	S	0	6	0
			887	544	163	174	6			

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		

- Molecule 4 is 2-[3-(2-HYDROXY-1,1-DIHYDROXYMETHYL-ETHYLAMINO)-PROPYL AMINO]-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: B3P) (formula: C₁₁H₂₆N₂O₆).

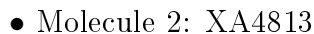
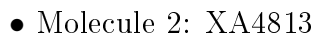


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			19	11	2	6		
4	B	1	Total	C	N	O	0	0
			19	11	2	6		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	210	Total	O	0	0
			210	210		
5	B	212	Total	O	0	0
			212	212		
5	D	96	Total	O	0	0
			96	96		
5	E	100	Total	O	0	0
			100	100		

- Molecule 1: BETA-SECRETASE 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.03 Å 61.76 Å 115.62 Å 90.00° 101.38° 90.00°	Depositor
Resolution (Å)	43.68 – 1.85 43.68 – 1.85	Depositor EDS
% Data completeness (in resolution range)	90.0 (43.68-1.85) 90.0 (43.68-1.85)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.22 (at 1.86 Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.181 , 0.222 0.183 , 0.224	Depositor DCC
R_{free} test set	3380 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	23.7	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8059	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 19.97 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 9.6362e-03.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: B3P, CL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.48	0/2889	0.62	0/3925
1	B	0.48	0/2901	0.65	0/3943
2	D	0.54	0/874	0.70	0/1182
2	E	0.49	0/905	0.63	0/1223
All	All	0.49	0/7569	0.64	0/10273

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	227	ASN	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2821	0	2734	24	0
1	B	2834	0	2742	18	0
2	D	860	0	835	6	0
2	E	887	0	869	15	0
3	A	1	0	0	0	0
4	A	19	0	26	2	0
4	B	19	0	26	1	0
5	A	210	0	0	2	0
5	B	212	0	0	3	0
5	D	96	0	0	3	0
5	E	100	0	0	13	0
All	All	8059	0	7232	61	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (61) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:246[A]:LYS:HE2	5:E:2083:HOH:O	1.71	0.88
2:E:178[A]:ARG:NH1	5:E:2022:HOH:O	2.14	0.80
2:E:236:ASN:CG	5:E:2080:HOH:O	2.19	0.79
2:E:246[A]:LYS:CE	5:E:2083:HOH:O	2.27	0.78
1:B:141[B]:THR:HG23	5:B:2091:HOH:O	1.82	0.78
2:D:271:SER:C	5:D:2016:HOH:O	2.22	0.78
2:E:178[A]:ARG:HD2	5:E:2022:HOH:O	1.84	0.77
2:E:178[B]:ARG:NE	5:E:2023:HOH:O	1.84	0.77
1:A:344:VAL:HG21	4:A:1399:B3P:H72	1.71	0.71
2:E:246[B]:LYS:HD3	5:E:2083:HOH:O	1.94	0.68
2:E:178[B]:ARG:NH2	5:E:2027:HOH:O	2.29	0.65
1:B:320:GLN:OE1	1:B:333[A]:ARG:NE	2.22	0.65
2:E:178[B]:ARG:NH2	5:E:2023:HOH:O	2.24	0.65
2:E:198:GLN:NE2	2:E:202:LYS:O	2.32	0.61
1:B:129:ILE:HD12	1:B:131:TRP:O	2.02	0.59
1:A:30:TYR:CE1	1:A:170:MET:HE2	2.38	0.59
1:B:26:SER:HA	1:B:351:GLU:OE2	2.03	0.58
1:A:25:ASP:HB2	5:A:2005:HOH:O	2.04	0.57
1:A:197:TYR:HA	1:A:370:PRO:HD3	1.86	0.56
1:A:111[B]:PHE:CZ	1:A:156:LEU:HD13	2.43	0.54
2:D:271:SER:HB2	5:D:2096:HOH:O	2.08	0.54
2:E:178[B]:ARG:CZ	5:E:2023:HOH:O	2.36	0.53
1:A:30:TYR:CE1	1:A:170:MET:CE	2.93	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:2112:HOH:O	2:D:257:SER:OG	2.19	0.50
1:B:61:HIS:HB3	1:B:64:ILE:HG12	1.94	0.49
1:A:158:THR:HG22	2:D:194:THR:HG21	1.94	0.49
1:B:259:GLU:HG3	1:B:263:ARG:HH11	1.78	0.49
1:A:126:LEU:O	1:A:129:ILE:HG12	2.13	0.48
1:A:265:SER:O	1:A:292:TYR:HE2	1.97	0.47
1:A:170:MET:CE	1:A:347:ALA:HB1	2.44	0.47
1:A:170:MET:HE1	1:A:347:ALA:HB1	1.96	0.47
1:A:139:TYR:CZ	1:A:209:GLU:HG2	2.49	0.47
1:B:58:GLY:HA3	1:B:117:THR:HG22	1.97	0.47
1:B:64:ILE:HD12	1:B:66[B]:THR:O	2.14	0.47
1:A:265:SER:O	1:A:292:TYR:CE2	2.69	0.46
1:A:34:MET:HE3	1:A:34:MET:HB3	1.68	0.46
1:A:246:LEU:HD22	1:A:335:GLY:HA2	1.96	0.46
1:B:109[A]:THR:HG22	5:B:2075:HOH:O	2.16	0.45
2:E:164:GLN:NE2	5:E:2007:HOH:O	2.48	0.45
1:B:35:LEU:HB3	1:B:40:PRO:HB2	1.99	0.44
1:A:26:SER:HB3	1:A:351:GLU:OE2	2.17	0.44
1:A:158:THR:CG2	2:D:194:THR:HG21	2.48	0.44
2:E:236:ASN:OD1	5:E:2080:HOH:O	2.20	0.44
1:A:322:MET:HE1	1:A:333[B]:ARG:NH1	2.32	0.44
1:A:170:MET:HE1	1:A:347:ALA:CB	2.48	0.43
1:B:35:LEU:HB2	1:B:102:THR:HB	2.00	0.43
4:A:1399:B3P:H31	4:A:1399:B3P:H61	1.66	0.42
1:A:34:MET:CE	1:A:54:PHE:CZ	3.02	0.42
1:B:89:GLN:HE21	1:B:89:GLN:HB2	1.70	0.42
1:A:111[A]:PHE:CE1	1:A:160:ALA:HB2	2.55	0.42
1:A:235:ALA:HA	1:A:236:ASP:HA	1.68	0.41
1:B:158[A]:THR:CG2	2:E:194:THR:HG21	2.50	0.41
1:B:163:PRO:HD2	1:B:190:GLY:O	2.20	0.41
1:B:288:THR:HB	1:B:289:PRO:HD2	2.02	0.41
2:D:202:LYS:HE2	5:D:2004:HOH:O	2.20	0.41
1:B:197:TYR:HA	1:B:370:PRO:HD3	2.03	0.41
1:B:78:SER:HB2	5:B:2052:HOH:O	2.20	0.40
1:A:34:MET:HE1	1:A:54:PHE:CZ	2.56	0.40
1:B:87:TYR:CZ	4:B:1398:B3P:H112	2.57	0.40
1:A:111[A]:PHE:CZ	1:A:160:ALA:HB2	2.57	0.40
2:E:236:ASN:CB	5:E:2080:HOH:O	2.65	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	354/386 (92%)	350 (99%)	4 (1%)	0	100	100
1	B	356/386 (92%)	351 (99%)	5 (1%)	0	100	100
2	D	114/122 (93%)	111 (97%)	3 (3%)	0	100	100
2	E	117/122 (96%)	116 (99%)	1 (1%)	0	100	100
All	All	941/1016 (93%)	928 (99%)	13 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	306/320 (96%)	303 (99%)	3 (1%)	76	69
1	B	308/320 (96%)	303 (98%)	5 (2%)	62	49
2	D	94/99 (95%)	93 (99%)	1 (1%)	73	65
2	E	97/99 (98%)	96 (99%)	1 (1%)	76	69
All	All	805/838 (96%)	795 (99%)	10 (1%)	71	62

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	210	TRP
1	A	284	THR

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Mol	Chain	Res	Type
1	A	393	SER
1	B	109[A]	THR
1	B	109[B]	THR
1	B	210	TRP
1	B	236	ASP
1	B	272	ASP
2	D	271	SER
2	E	221	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	89	GLN
1	B	169	GLN
1	B	341	ASN
2	D	164	GLN
2	E	164	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	B3P	A	1399	-	18,18,18	0.51	0	21,23,23	1.49	2 (9%)
4	B3P	B	1398	-	18,18,18	0.54	0	21,23,23	2.24	4 (19%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	B3P	A	1399	-	-	16/28/28/28	-
4	B3P	B	1398	-	-	13/28/28/28	-

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1398	B3P	C3-N1-C4	7.56	126.80	116.08
4	B	1398	B3P	C2-N2-C8	5.68	124.13	116.08
4	A	1399	B3P	C3-N1-C4	4.42	122.34	116.08
4	A	1399	B3P	C2-N2-C8	3.93	121.65	116.08
4	B	1398	B3P	C6-C4-C5	-2.23	105.34	110.04
4	B	1398	B3P	O2-C10-C8	-2.22	107.14	111.63

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1399	B3P	C5-C4-N1-C3
4	A	1399	B3P	C6-C4-N1-C3
4	A	1399	B3P	N1-C4-C5-O4
4	A	1399	B3P	C6-C4-C5-O4
4	A	1399	B3P	C7-C4-C5-O4
4	A	1399	B3P	N1-C4-C6-O5
4	A	1399	B3P	C5-C4-C6-O5
4	A	1399	B3P	C7-C4-C6-O5
4	A	1399	B3P	N1-C4-C7-O6
4	A	1399	B3P	C5-C4-C7-O6

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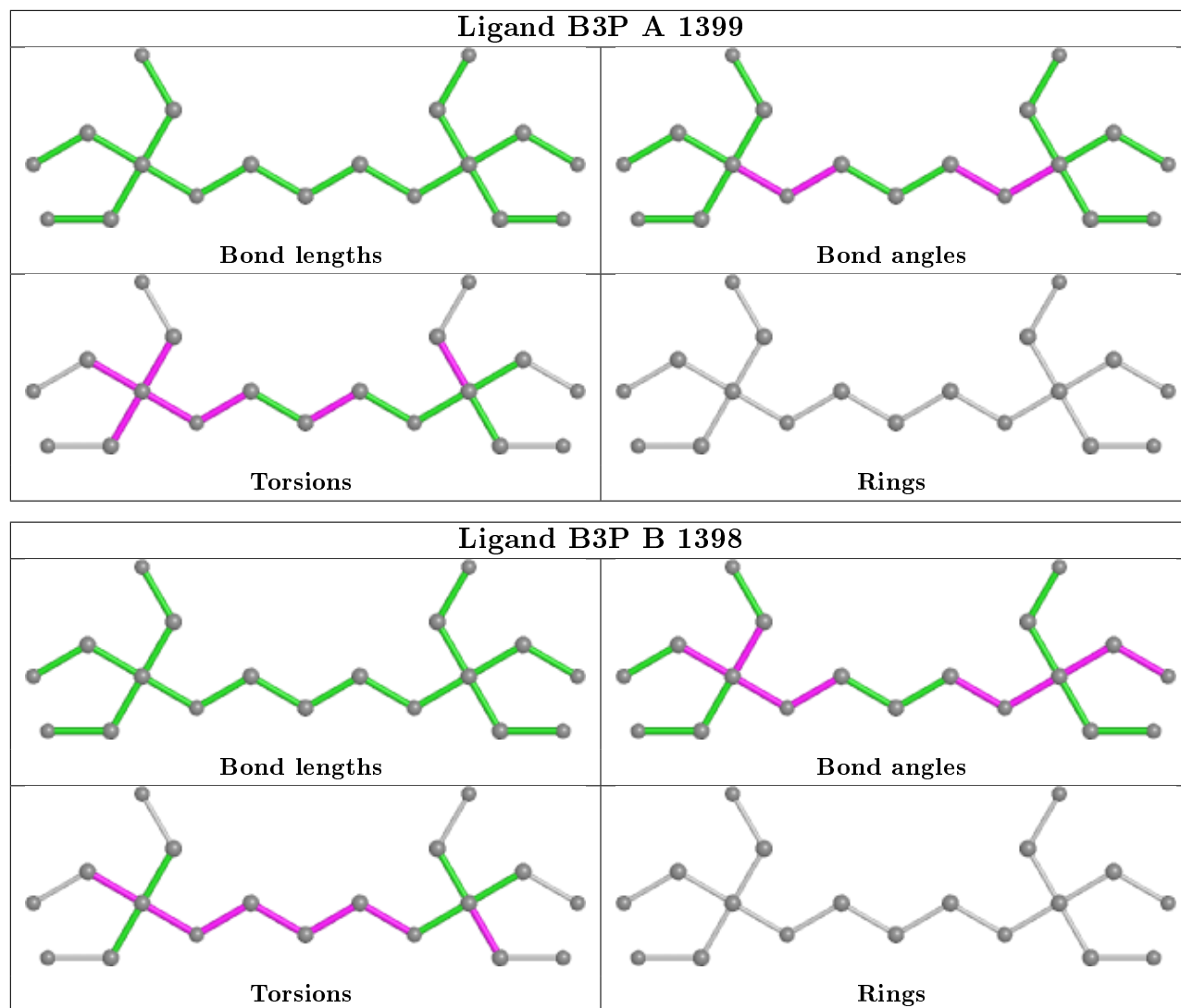
Mol	Chain	Res	Type	Atoms
4	A	1399	B3P	C6-C4-C7-O6
4	B	1398	B3P	C1-C3-N1-C4
4	B	1398	B3P	C5-C4-N1-C3
4	B	1398	B3P	C6-C4-N1-C3
4	B	1398	B3P	C7-C4-N1-C3
4	B	1398	B3P	N2-C8-C9-O1
4	B	1398	B3P	C11-C8-C9-O1
4	A	1399	B3P	C3-C1-C2-N2
4	B	1398	B3P	C3-C1-C2-N2
4	B	1398	B3P	C10-C8-C9-O1
4	B	1398	B3P	C2-C1-C3-N1
4	B	1398	B3P	N1-C4-C6-O5
4	B	1398	B3P	C7-C4-C6-O5
4	B	1398	B3P	C1-C2-N2-C8
4	A	1399	B3P	C7-C4-N1-C3
4	A	1399	B3P	C1-C3-N1-C4
4	A	1399	B3P	O3-C11-C8-C10
4	B	1398	B3P	C5-C4-C6-O5
4	A	1399	B3P	O3-C11-C8-N2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1399	B3P	2	0
4	B	1398	B3P	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
1	A	358/386 (92%)	-0.16	21 (5%)	22	22	18, 28, 59, 78	20 (5%)
1	B	357/386 (92%)	-0.24	16 (4%)	33	32	17, 27, 54, 76	20 (5%)
2	D	112/122 (91%)	-0.60	0	100	100	16, 23, 35, 57	4 (3%)
2	E	113/122 (92%)	-0.52	0	100	100	16, 24, 41, 53	2 (1%)
All	All	940/1016 (92%)	-0.29	37 (3%)	39	38	16, 26, 55, 78	46 (4%)

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	292	TYR	7.4
1	A	283	TRP	4.9
1	A	264	ALA	4.1
1	A	282	CYS	4.0
1	B	283	TRP	3.8
1	B	227	ASN	3.7
1	A	331	CYS	3.6
1	A	278	SER	3.4
1	A	276	THR	3.2
1	A	275	TRP	3.2
1	B	262	ALA	3.2
1	B	289	PRO	3.2
1	B	278	SER	2.9
1	B	27	GLY	2.9
1	B	63	TYR	2.8
1	B	88	THR	2.8
1	A	281	ALA	2.7
1	B	282	CYS	2.6
1	B	331	CYS	2.5
1	B	330	GLU	2.4
1	B	275	TRP	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	28	ARG	2.4
1	B	292	TYR	2.3
1	A	333[A]	ARG	2.3
1	A	88	THR	2.3
1	A	265	SER	2.3
1	A	280	LEU	2.2
1	A	89	GLN	2.2
1	A	293	PHE	2.2
1	B	89	GLN	2.2
1	A	26	SER	2.1
1	B	28	ARG	2.1
1	A	397	PRO	2.1
1	A	262	ALA	2.1
1	A	273	GLY	2.1
1	B	126	LEU	2.0
1	A	332	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

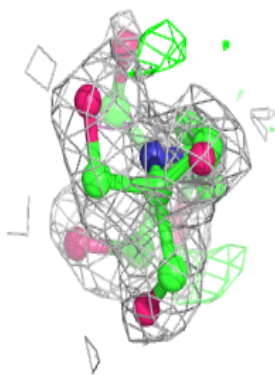
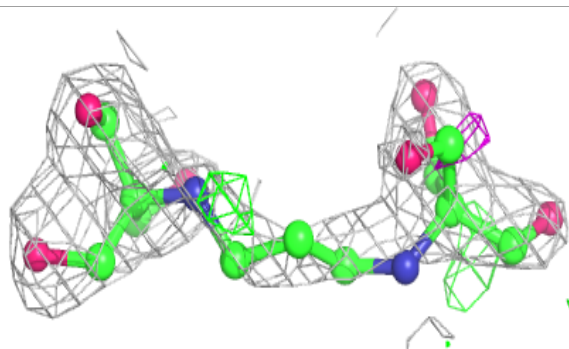
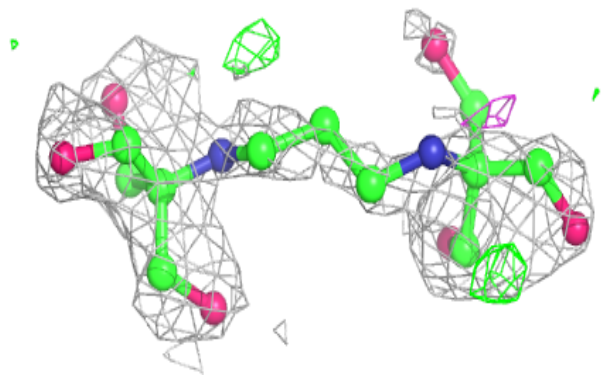
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	B3P	A	1399	19/19	0.72	0.22	41,45,47,47	19
4	B3P	B	1398	19/19	0.84	0.18	42,50,55,66	0
3	CL	A	1398	1/1	1.00	0.05	21,21,21,21	0

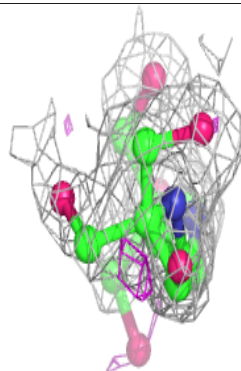
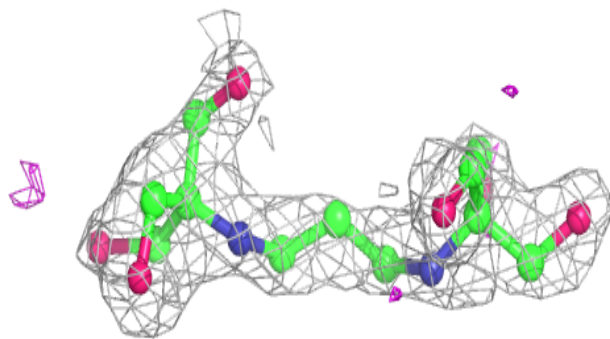
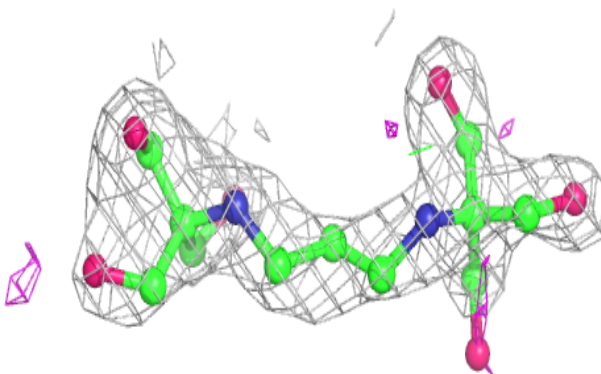
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around B3P A 1399:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around B3P B 1398:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.